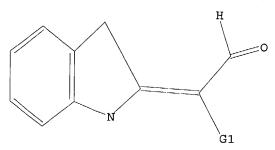
Uploading C:\Program Files\Stnexp\Queries\10055664.str

L1STRUCTURE UPLOADED

=> d

L1 HAS NO ANSWERS

L1



G1 CH,O

Structure attributes must be viewed using STN Express query preparation.

=> s 11

SAMPLE SEARCH INITIATED 10:15:33 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED -59 TO ITERATE

100.0% PROCESSED 59 ITERATIONS

0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS:

720 TO 1640 PROJECTED ANSWERS: O TO

L20 SEA SSS SAM L1

=> s l1 full

FULL SEARCH INITIATED 10:15:37 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 1283 TO ITERATE

100.0% PROCESSED 1283 ITERATIONS

2 ANSWERS

SEARCH TIME: 00.00.01

1.3 2 SEA SSS FUL L1

=> d 1-2

ANSWER 1 OF 2 REGISTRY COPYRIGHT 2004 ACS on STN L3

RN124244-70-8 REGISTRY

CNPropanal, 2-(1,3-dihydro-1,3,3-trimethyl-2H-indol-2-ylidene)-, (E)- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MFC14 H17 N O

SR CA

LCSTN Files: BEILSTEIN*, CA, CAPLUS

(*File contains numerically searchable property data)

DT.CA CAplus document type: Journal

RL.NP Roles from non-patents: PREP (Preparation); RACT (Reactant or reagent)

Double bond geometry as shown.

Moderna

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 1 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- L3 ANSWER 2 OF 2 REGISTRY COPYRIGHT 2004 ACS on STN

RN 42909-52-4 REGISTRY

CN Propanal, 2-(1,3-dihydro-1,3,3-trimethyl-2H-indol-2-ylidene)- (9CI) (CA INDEX NAME)

OTHER NAMES:

CN 1,3,3-Trimethyl-2-(α -formylethylidene)indoline

CN 1,3,3-Trimethyl-2-(1-formylethylidene)indoline

FS 3D CONCORD

MF C14 H17 N O

LC STN Files: BEILSTEIN*, CA, CAPLUS

(*File contains numerically searchable property data)

DT.CA CAplus document type: Patent

RL.P Roles from patents: PREP (Preparation); USES (Uses)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 3 REFERENCES IN FILE CA (1907 TO DATE)
- 3 REFERENCES IN FILE CAPLUS (1907 TO DATE)

>>> FOR SEARCHING PREPARATIONS SEE HELP PRE <<< ****************** * PLEASE NOTE THAT THERE ARE NO FORMATS FREE OF COST. * SET NOTICE FEATURE: THE COST ESTIMATES CALCULATED FOR SET NOTICE * ARE BASED ON THE HIGHEST PRICE CATEGORY. THEREFORE; THESE * ESTIMATES MAY NOT REFLECT THE ACTUAL COSTS. * FOR PRICE INFORMATION SEE HELP COST ********************** => s l1 full FULL SEARCH INITIATED 10:16:14 FILE 'BEILSTEIN' FULL SCREEN SEARCH COMPLETED - 651 TO ITERATE 100.0% PROCESSED 651 ITERATIONS 1 ANSWERS SEARCH TIME: 00.00.10 L4 1 SEA SSS FUL L1 => d ide ANSWER 1 OF 1 BEILSTEIN COPYRIGHT 2004 BEILSTEIN MDL on STN L4Beilstein Records (BRN): 3545115 Beilstein Pref. RN (BPR): 124244-70-8 CAS Reg. No. (RN): 42909-52-4, 124244-70-8 Chemical Name (CN): 2-(1,3,3-Trimethyl-2indolinyliden)propanal Autonom Name (AUN): ylidene)-propionaldehyde Molec. Formula (MF): C14 H17 N O Molecular Weight (MW): 215.29 Lawson Number (LN): 25479, 2817

indolinyliden)propanal
2-(1,3,3-trimethyl-1,3-dihydro-indol-2ylidene)-propionaldehyde

Molec. Formula (MF):

Molecular Weight (MW):

Lawson Number (LN):

File Segment (FS):

Constitution ID (CONSID):

Tautomer ID (TAUTID):

Beilstein Citation (BSO):

Entry Date (DUPD):

indolinyliden)propanal

2-(1,3,3-trimethyl-1,3-dihydro-indol-2ylidene)-propionaldehyde

C14 H17 N O

25479, 2817

Stereo compound
heterocyclic

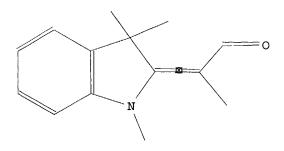
3137287

3361259

Beilstein Citation (BSO):
6-21

Entry Date (DED):
1991/10/23

Update Date (DUPD):
1991/11/01



Field Availability:

| Code | Name | Occurrence | | |
|---|-------------------|------------|--|--|
| ======================================= | | | | |
| BRN | Beilstein Records | | | |

| BPR | Beilstein Preferred RN | 1 |
|--------|----------------------------|---|
| RN | CAS Registry Number | 2 |
| CN | Chemical Name | 1 |
| AUN | Autonomname | 1 |
| MF | Molecular Formula | 1 |
| FW | Formular Weight | 1 |
| LN | Lawson Number | 2 |
| FS | File Segment | ī |
| CTYPE | Compound Type | 1 |
| CONSID | Constitution ID | 1 |
| TAUTID | Tautomer ID | 1 |
| BSO | Beilstein Citation | 1 |
| ED | Entry Date | 1 |
| UPD | Update Date | 1 |
| MP | Melting Point | 1 |
| NMR | Nuclear Magnetic Resonance | 3 |
| UVS | UV and Visible Spectrum | 1 |
| | F | - |

This substance also occurs in Reaction Documents:

| Code | Name | Occurrence |
|---------|--|-------------|
| ======= | ====================================== | =========== |
| RX | Reaction Documents | 3 |
| RXREA | Substance is Reaction Reactant | 2 |
| RXPRO | Substance is Reaction Product | 1 |

Connection closed by remote host